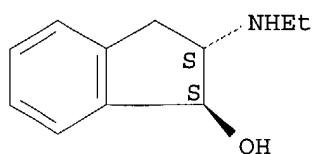


CAS ONLINE PRINTOUT

=> d scan

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2-(ethylamino)-2,3-dihydro-, (1R,2R)-rel- (9CI)
MF C11 H15 N O
CI COM

Relative stereochemistry.

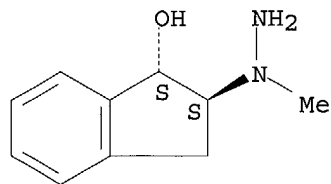


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):50

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1S,2S)- (9CI)
MF C10 H14 N2 O
CI COM

Absolute stereochemistry. Rotation (+).



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

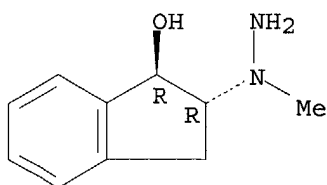
L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2R)-, (2Z)-2-butenedioate (1:1) (salt) (9CI)
MF C10 H14 N2 O . C4 H4 O4

CM 1

Absolute stereochemistry.

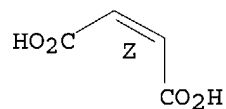
compounds disclosed in specification

CAS ONLINE PRINTOUT



CM 2

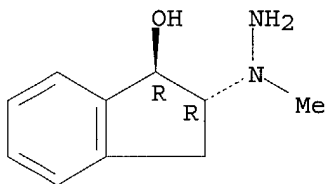
Double bond geometry as shown.



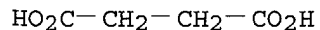
L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN Butanedioic acid, compd. with (1R,2R)-2,3-dihydro-2-(1-methylhydrazino)-1H-inden-1-ol (1:2) (9CI)
MF C10 H14 N2 O . 1/2 C4 H6 O4

CM 1

Absolute stereochemistry.



CM 2



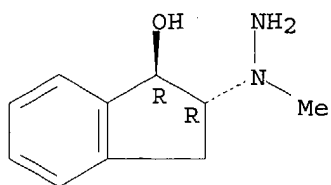
L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2R)-, (2R,3R)-2,3-dihydroxybutanedioate (2:1) (salt) (9CI)
MF C10 H14 N2 O . 1/2 C4 H6 O6

CM 1

Absolute stereochemistry.

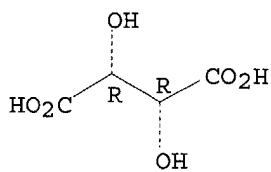
compounds disclosed in specification

CAS ONLINE PRINTOUT



CM 2

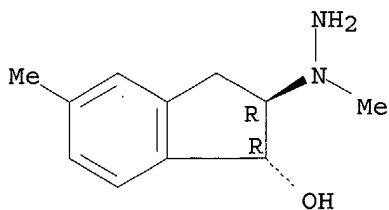
Absolute stereochemistry.



L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
 IN 1H-Inden-1-ol, 2,3-dihydro-5-methyl-2-(1-methylhydrazino)-, (1R,2R)-rel-,
 (2Z)-2-butenedioate (1:1) (salt) (9CI)
 MF C11 H16 N2 O . C4 H4 O4

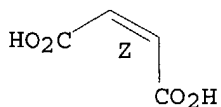
CM 1

Relative stereochemistry.



CM 2

Double bond geometry as shown.

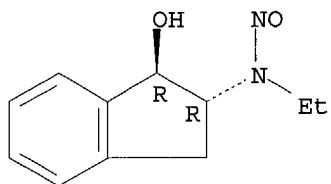


L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
 IN 1H-Inden-1-ol, 2-(ethylnitrosoamino)-2,3-dihydro-, (1R,2R)-rel- (9CI)
 MF C11 H14 N2 O2

Relative stereochemistry.

compounds disclosed in specification

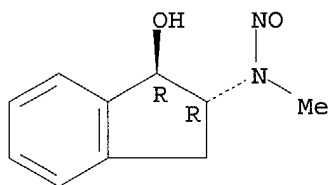
CAS ONLINE PRINTOUT



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(methylnitrosoamino)-, (1R,2R)-rel- (9CI)
MF C10 H12 N2 O2

Relative stereochemistry.

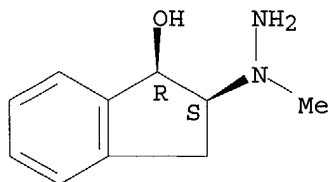


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2S)-,
(2Z)-2-butenedioate (1:1) (salt) (9CI)
MF C10 H14 N2 O . C4 H4 O4

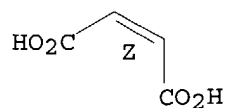
CM 1

Absolute stereochemistry.



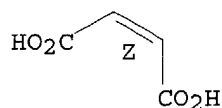
CM 2

Double bond geometry as shown.



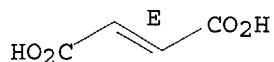
compounds disclosed in specification

CAS ONLINE PRINTOUT



L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 2-Butenedioic acid (2E)- (9CI)
MF C4 H4 O4
CI COM

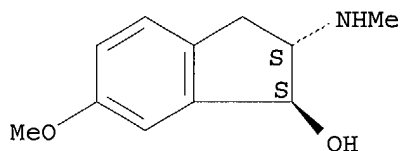
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-6-methoxy-2-(methylamino)-, (1R,2R)-rel- (9CI)
MF C11 H15 N O2
CI COM

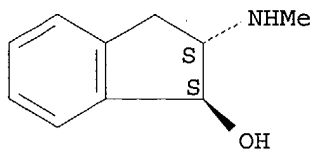
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(methylamino)-, (1S-trans)- (9CI)
MF C10 H13 N O

Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2R)-rel-,

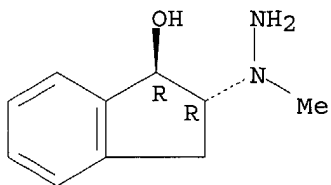
compounds disclosed in specification

CAS ONLINE PRINTOUT

(2Z)-2-butenedioate (1:1) (salt) (9CI)
MF C10 H14 N2 O . C4 H4 O4

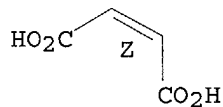
CM 1

Relative stereochemistry.



CM 2

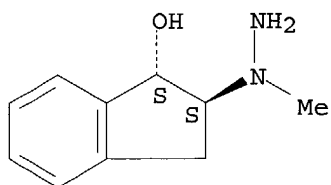
Double bond geometry as shown.



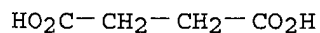
L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN Butanedioic acid, compd. with (1S,2S)-2,3-dihydro-2-(1-methylhydrazino)-1H-inden-1-ol (1:2) (9CI)
MF C10 H14 N2 O . 1/2 C4 H6 O4

CM 1

Absolute stereochemistry. Rotation (+).



CM 2



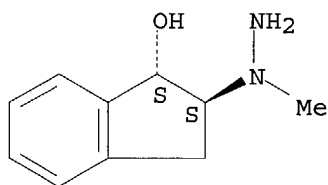
L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1S,2S)-, (2R,3R)-2,3-dihydroxybutanedioate (2:1) (salt) (9CI)
MF C10 H14 N2 O . 1/2 C4 H6 O6

compounds disclosed in specification

CAS ONLINE PRINTOUT

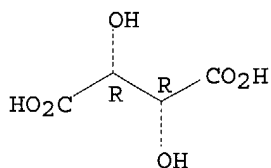
CM 1

Absolute stereochemistry. Rotation (+).



CM 2

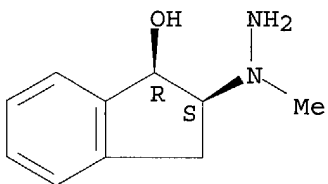
Absolute stereochemistry.



L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
 IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2S)-rel-,
 (2Z)-2-butenedioate (1:1) (salt) (9CI)
 MF C10 H14 N2 O . C4 H4 O4

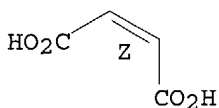
CM 1

Relative stereochemistry.



CM 2

Double bond geometry as shown.



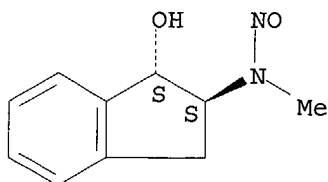
L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
 IN 1H-Inden-1-ol, 2,3-dihydro-2-(methylnitrosoamino)-, (1S,2S)- (9CI)

compounds disclosed in specification

CAS ONLINE PRINTOUT

MF C10 H12 N2 O2

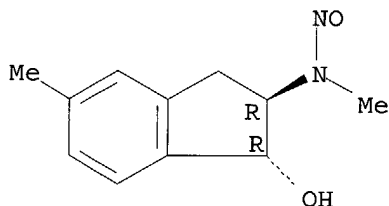
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-5-methyl-2-(methylnitrosoamino)-, (1R,2R)-rel- (9CI)
MF C11 H14 N2 O2

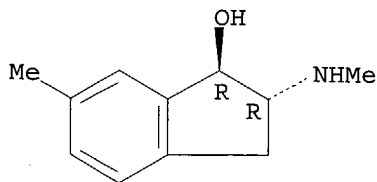
Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-6-methyl-2-(methylamino)-, (1R,2R)-rel- (9CI)
MF C11 H15 N O

Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

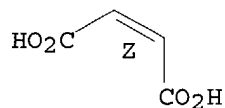
L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN

compounds disclosed in specification

CAS ONLINE PRINTOUT

IN 2-Butenedioic acid (2Z)- (9CI)
MF C4 H4 O4
CI COM

Double bond geometry as shown.



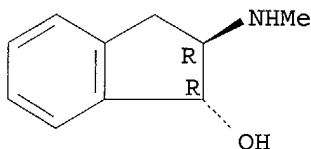
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN Oxidase, monoamine (9CI)
MF Unspecified
CI MAN

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(methylamino)-, (1R-trans)- (9CI)
MF C10 H13 N O

Absolute stereochemistry.

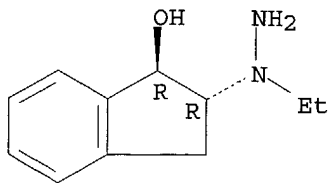


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2-(1-ethylhydrazino)-2,3-dihydro-, (1R,2R)-rel-,
(2Z)-2-butenedioate (1:1) (salt) (9CI)
MF C11 H16 N2 O . C4 H4 O4

CM 1

Relative stereochemistry.

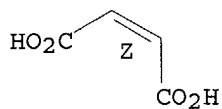


compounds disclosed in specification

CAS ONLINE PRINTOUT

CM 2

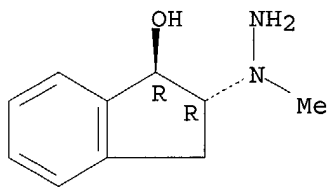
Double bond geometry as shown.



L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2R)-,
(2E)-2-butenedioate (2:1) (salt) (9CI)
MF C10 H14 N2 O . 1/2 C4 H4 O4

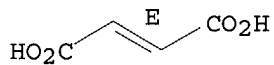
CM 1

Absolute stereochemistry.



CM 2

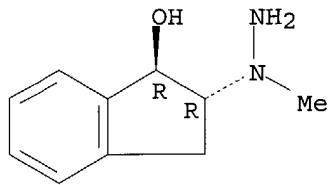
Double bond geometry as shown.



L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2R)-,
(2S,3S)-2,3-dihydroxybutanedioate (2:1) (salt) (9CI)
MF C10 H14 N2 O . 1/2 C4 H6 O6

CM 1

Absolute stereochemistry.

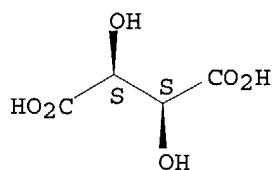


CM 2

compounds disclosed in specification

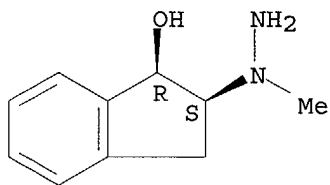
CAS ONLINE PRINTOUT

Absolute stereochemistry.



L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1R,2S)- (9CI)
MF C10 H14 N2 O
CI COM

Absolute stereochemistry.

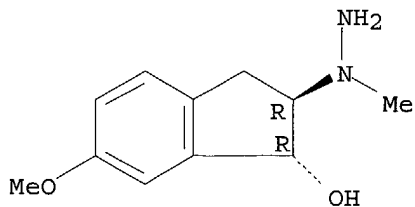


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-6-methoxy-2-(1-methylhydrazino)-, (1R,2R)-rel-,
(2Z)-2-butenedioate (1:1) (salt) (9CI)
MF C11 H16 N2 O2 . C4 H4 O4

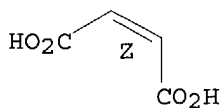
CM 1

Relative stereochemistry.



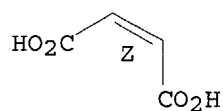
CM 2

Double bond geometry as shown.



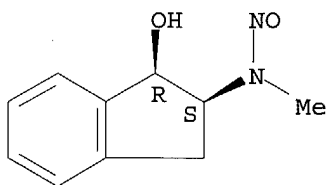
compounds disclosed in specification

CAS ONLINE PRINTOUT



L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(methylnitrosoamino)-, (1R,2S)- (9CI)
MF C10 H12 N2 O2

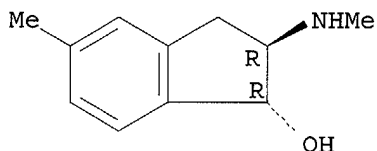
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

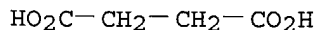
L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-5-methyl-2-(methylamino)-, (1R,2R)-rel- (9CI)
MF C11 H15 N O

Relative stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN Butanedioic acid (9CI)
MF C4 H6 O4
CI COM



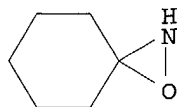
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1-Oxa-2-azaspiro[2.5]octane (7CI, 8CI, 9CI)
MF C6 H11 N O

compounds disclosed in specification

CAS ONLINE PRINTOUT

CI RPS



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

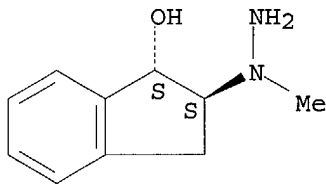
L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN VAP 1 (9CI)
MF Unspecified
CI MAN

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1S,2S)-, (2Z)-2-butenedioate (1:1) (salt) (9CI)
MF C10 H14 N2 O . C4 H4 O4

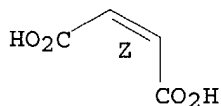
CM 1

Absolute stereochemistry. Rotation (+).



CM 2

Double bond geometry as shown.



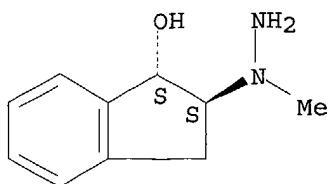
L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1S,2S)-, (2E)-2-butenedioate (2:1) (salt) (9CI)
MF C10 H14 N2 O . 1/2 C4 H4 O4

CM 1

Absolute stereochemistry. Rotation (+).

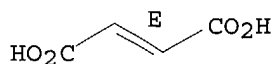
compounds disclosed in specification

CAS ONLINE PRINTOUT



CM 2

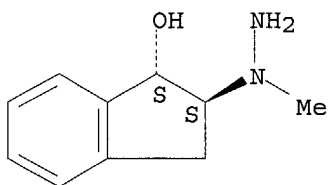
Double bond geometry as shown.



L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
 IN 1H-Inden-1-ol, 2,3-dihydro-2-(1-methylhydrazino)-, (1S,2S)-,
 (2S,3S)-2,3-dihydroxybutanedioate (2:1) (salt) (9CI)
 MF C10 H14 N2 O . 1/2 C4 H6 O6

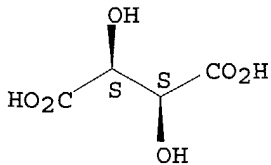
CM 1

Absolute stereochemistry. Rotation (+).



CM 2

Absolute stereochemistry.

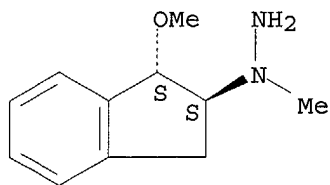


L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
 IN Hydrazine, 1-[(1S,2S)-2,3-dihydro-1-methoxy-1H-inden-2-yl]-1-methyl-,
 (2Z)-2-butenedioate (1:1) (9CI)
 MF C11 H16 N2 O . C4 H4 O4

CM 1

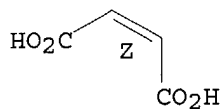
Absolute stereochemistry.

compounds disclosed in specification



CM 2

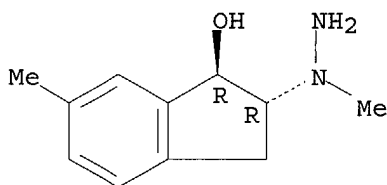
Double bond geometry as shown.



L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-6-methyl-2-(1-methylhydrazino)-, (1R,2R)-rel-,
(2Z)-2-butenedioate (1:1) (salt) (9CI)
MF C11 H16 N2 O . C4 H4 O4

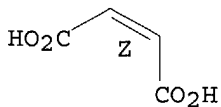
CM 1

Relative stereochemistry.



CM 2

Double bond geometry as shown.

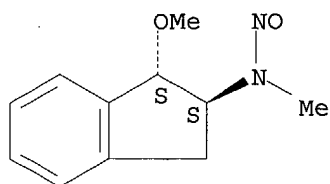


L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-2-amine, 2,3-dihydro-1-methoxy-N-methyl-N-nitroso-, (1S,2S)-
(9CI)
MF C11 H14 N2 O2

Absolute stereochemistry.

compounds disclosed in specification

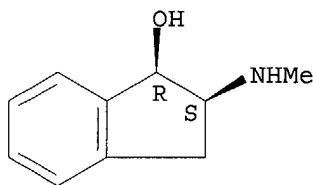
CAS ONLINE PRINTOUT



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN 1H-Inden-1-ol, 2,3-dihydro-2-(methylanilino)-, (1R,2S)- (9CI)
MF C10 H13 N O

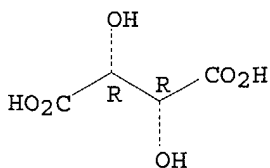
Absolute stereochemistry.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN Butanedioic acid, 2,3-dihydroxy-, (2R,3R)- (9CI)
MF C4 H6 O6
CI COM

Absolute stereochemistry.



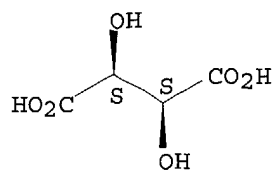
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L7 41 ANSWERS REGISTRY COPYRIGHT 2003 ACS on STN
IN Butanedioic acid, 2,3-dihydroxy-, (2S,3S)- (9CI)
MF C4 H6 O6
CI COM

Absolute stereochemistry.

compounds disclosed in specification

CAS ONLINE PRINTOUT



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=>

compounds disclosed in specification

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RL: SPN (Synthetic preparation); PREP (Preparation)

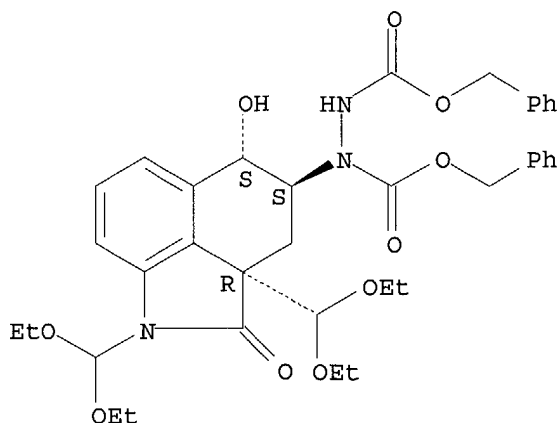
(synthesis, pharmacol. investigation and computational studies on a tricyclic ergoline analog with selective dopamine autoreceptor activity)

09/902,789

RN 171006-44-3 CAPLUS

CN 1,2-Hydrazinedicarboxylic acid, 1-[1,2a-bis(diethoxymethyl)-1,2,2a,3,4,5-hexahydro-5-hydroxy-2-oxobenz[cd]indol-4-yl]-, bis(phenylmethyl) ester, (2a.alpha.,4.beta.,5.alpha.)- (9CI) (CA INDEX NAME)

Relative stereochemistry.



L14 ANSWER 4 OF 7 CAPLUS COPYRIGHT 2003 ACS on STN

AN 1992:426035 CAPLUS

DN 117:26035

TI Synthesis of amines and amino alcohols by electrophilic amination and highly stereoselective reduction

AU Gmeiner, Peter; Bollinger, Bernd

CS Inst. Pharm. Lebensmittelchem., Univ. Muenchen, Munich, 8000/2, Germany

SO Liebigs Annalen der Chemie (1992), (3), 273-8

CODEN: LACHDL; ISSN: 0170-2041

DT Journal

LA English

OS CASREACT 117:26035

IT 138206-94-7P 138206-95-8P 138408-15-8P

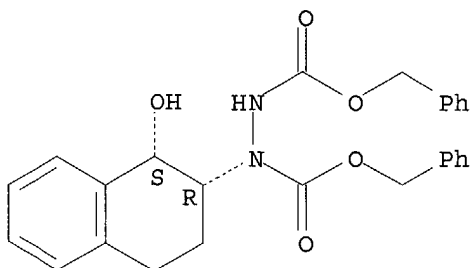
138408-16-9P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent) (prepn. and hydrogenolysis of)

RN 138206-94-7 CAPLUS

CN 1,2-Hydrazinedicarboxylic acid, 1-(1,2,3,4-tetrahydro-1-hydroxy-2-naphthalenyl)-, bis(phenylmethyl) ester, cis- (9CI) (CA INDEX NAME)

Relative stereochemistry.



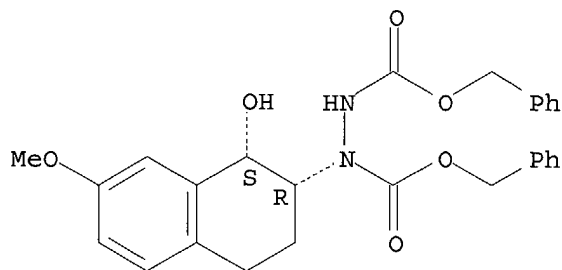
RN 138206-95-8 CAPLUS

CN 1,2-Hydrazinedicarboxylic acid, 1-(1,2,3,4-tetrahydro-1-hydroxy-7-methoxy-

09/902,789

2-naphthalenyl)-, bis(phenylmethyl) ester, cis- (9CI) (CA INDEX NAME)

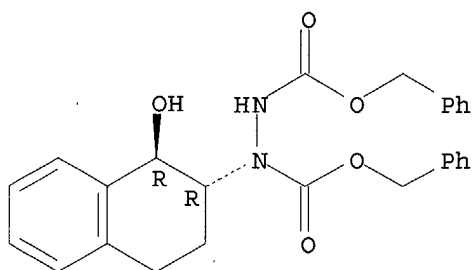
Relative stereochemistry.



RN 138408-15-8 CAPLUS

CN 1,2-Hydrazinedicarboxylic acid, 1-(1,2,3,4-tetrahydro-1-hydroxy-2-naphthalenyl)-, bis(phenylmethyl) ester, trans- (9CI) (CA INDEX NAME)

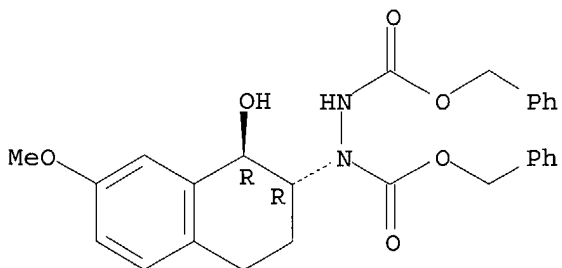
Relative stereochemistry.



RN 138408-16-9 CAPLUS

CN 1,2-Hydrazinedicarboxylic acid, 1-(1,2,3,4-tetrahydro-1-hydroxy-7-methoxy-2-naphthalenyl)-, bis(phenylmethyl) ester, trans- (9CI) (CA INDEX NAME)

Relative stereochemistry.



IT 138408-13-6P 138408-14-7P 138408-17-0P
138408-18-1P

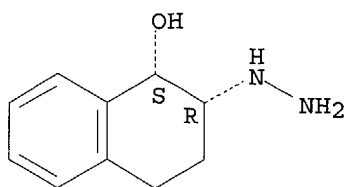
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT
(Reactant or reagent)
(prepn. and nitrogen-nitrogen bond cleavage of)

RN 138408-13-6 CAPLUS

CN 1-Naphthalenol, 2-hydrazino-1,2,3,4-tetrahydro-, monohydrochloride, cis-
(9CI) (CA INDEX NAME)

Relative stereochemistry.

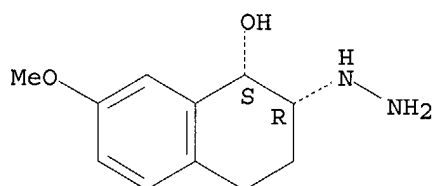
09/902,789



● HCl

RN 138408-14-7 CAPLUS
CN 1-Naphthalenol, 2-hydrazino-1,2,3,4-tetrahydro-7-methoxy-,
monohydrochloride, cis- (9CI) (CA INDEX NAME)

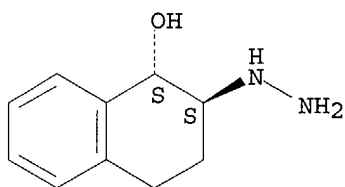
Relative stereochemistry.



● HCl

RN 138408-17-0 CAPLUS
CN 1-Naphthalenol, 2-hydrazino-1,2,3,4-tetrahydro-, monohydrochloride, trans-
(9CI) (CA INDEX NAME)

Relative stereochemistry.

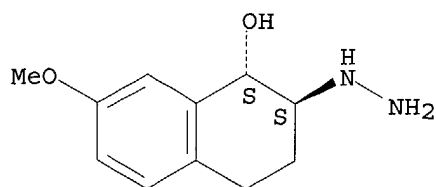


● HCl

RN 138408-18-1 CAPLUS
CN 1-Naphthalenol, 2-hydrazino-1,2,3,4-tetrahydro-7-methoxy-,
monohydrochloride, trans- (9CI) (CA INDEX NAME)

Relative stereochemistry.

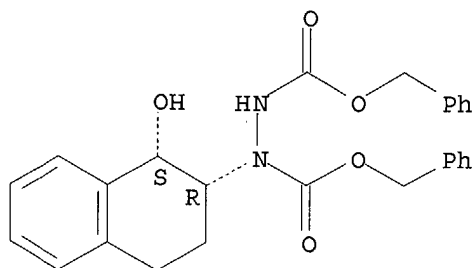
09/902,789



● HCl

L14 ANSWER 5 OF 7 CAPLUS COPYRIGHT 2003 ACS on STN
AN 1992:41818 CAPLUS
DN 116:41818
TI Efficient methodology for the preparation of .beta.-aminotetralin
derivatives via electrophilic amination
AU Gmeiner, Peter; Bollinger, Bernd
CS Inst. Pharm. Lebensmittelchem., Ludwig-Maximilians-Univ., Munich, 8000/2,
Germany
SO Tetrahedron Letters (1991), 32(42), 5927-30
CODEN: TELEAY; ISSN: 0040-4039
DT Journal
LA English
OS CASREACT 116:41818
IT **138206-94-7P 138206-95-8P**
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT
(Reactant or reagent)
(prepn. and transesterification-cyclization of)
RN 138206-94-7 CAPLUS
CN 1,2-Hydrazinedicarboxylic acid, 1-(1,2,3,4-tetrahydro-1-hydroxy-2-
naphthalenyl)-, bis(phenylmethyl) ester, cis- (9CI) (CA INDEX NAME)

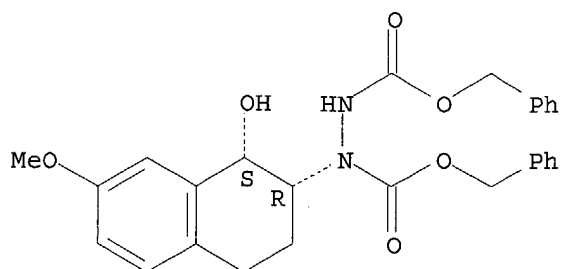
Relative stereochemistry.



RN 138206-95-8 CAPLUS
CN 1,2-Hydrazinedicarboxylic acid, 1-(1,2,3,4-tetrahydro-1-hydroxy-7-methoxy-
2-naphthalenyl)-, bis(phenylmethyl) ester, cis- (9CI) (CA INDEX NAME)

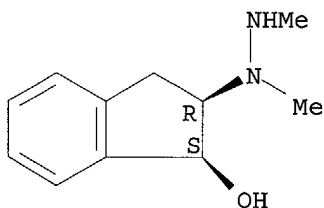
Relative stereochemistry.

09/902,789



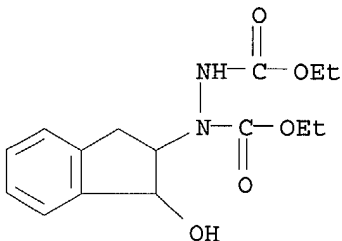
L14 ANSWER 6 OF 7 CAPLUS COPYRIGHT 2003 ACS on STN
AN 1970:121454 CAPLUS
DN 72:121454
TI Photochemical and thermal 1,2- and 1,4-cycloaddition reactions of
azodicarbonyl compounds with monoolefins
AU Koerner von Gustorf, Ernst; White, Danny V.; Kim, Bongsub; Hess, Dieter;
Leitich, Johannes
CS Abt. Strahlenchem., Max Planck Inst. Kohlenforsch., Muelheim, Fed. Rep.
Ger.
SO Journal of Organic Chemistry (1970), 35(4), 1155-65
CODEN: JOCEAH; ISSN: 0022-3263
DT Journal
LA English
IT 23358-19-2P 23358-23-8P
RL: SPN (Synthetic preparation); PREP (Preparation)
(prepn. of)
RN 23358-19-2 CAPLUS
CN 1-Indanol, 2-(1,2-dimethylhydrazino)-, cis- (8CI) (CA INDEX NAME)

Relative stereochemistry.



*Proviso
ent*

RN 23358-23-8 CAPLUS
CN Bicarbamidic acid, (1-hydroxy-2-indanyl)-, diethyl ester (8CI) (CA INDEX NAME)



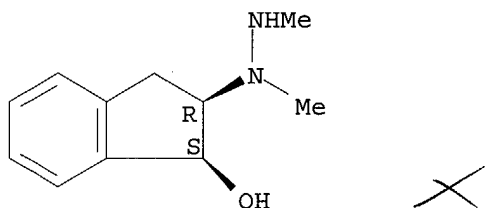
X

L14 ANSWER 7 OF 7 CAPLUS COPYRIGHT 2003 ACS on STN
AN 1970:121451 CAPLUS

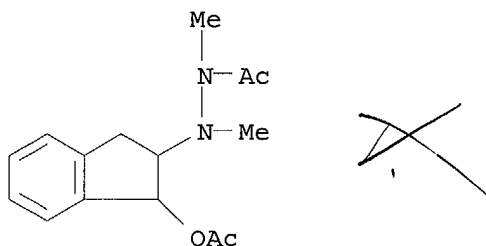
09/902,789

DN 72:121451
TI Azodiformate adduct of indene and the stereochemistry of some
1,2-disubstituted indans
AU Huebner, Charles F.; Donoghue, Ellen M.; Novak, Carol J.; Dorfman, Louis;
Wenkert, Ernest
CS Chem. Res. Div., CIBA Pharm. Co., Summit, NJ, USA
SO Journal of Organic Chemistry (1970), 35(4), 1149-54
CODEN: JOCEAH; ISSN: 0022-3263
DT Journal
LA English
IT **23358-19-2P 23359-96-8P**
RL: SPN (Synthetic preparation); PREP (Preparation)
(prepn. of)
RN 23358-19-2 CAPLUS
CN 1-Indanol, 2-(1,2-dimethylhydrazino)-, cis- (8CI) (CA INDEX NAME)

Relative stereochemistry.



RN 23359-96-8 CAPLUS
CN Acetic acid, 2-(1-hydroxy-2-indanyl)-1,2-dimethylhydrazide acetate (ester)
(8CI) (CA INDEX NAME)



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